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Information for Contributorsv
UPAC-NIST Solubility Data Series. 79. Alkali and Alkaline Earth Metal Pseudohalides
This volume presents solubility data of azides, cyanides, cyanates, and thiocyanates of alkali metals, alkaline earth metals, and ammonium. Covered are binary and ternary systems in all solvents. The literature has been covered up to the middle of 2001, and there was a great effort to have the literature survey as complete as cossible.
NIST Recommended Rest Frequencies for Observed Interstellar Molecular Microwave Transitions—2002 Revision
Critically evaluated transition frequencies for the molecular transitions detected in interstellar and circumstellar clouds are presented. The tabulated transitions are recommended for reference in future astronomical observations in the microwave and millimeter wavelength regions. The transition frequencies have been selected through a critical examination and analysis of the laboratory spectral data obtained from the literature. The information tabulated includes the species identity, transition frequency, uncertainty, and quantum state labels. For convenience, representative line antenna temperatures are listed for a typical astronomical source for each transition, and the references are cited for the laboratory and astronomical literature that have been employed.
Revised and Updated Thermochemical Properties of the Gases Mercapto (HS), Disulfur Monoxide ( $S_2O$ ), Thiazyl (NS), and Thioxophosphino (PS)
Computations were done to correct erroneous data tables given in the 4th edition of the NIST-JANAF Ther- mochemical Tables. Updated enthalpies of formation were included to compute the thermochemical tables for four ideal gases.
Correlation for the Second Virial Coefficient of Water
A new correlation has been developed to represent the second virial coefficient of water ( $H_2O$ ) as a function of temperature. The new correlation agrees well with the experimental data deemed to be reliable, and at high temperatures is a significant improvement over the best previous formulation.
Structure and Vibrations of Lanthanide Trihalides. An Assessment of Experimental and Theoretical Data

An assessment of experimental and theoretical data on the structure and molecular vibrations of all the  $LnX_3$  lanthanide trihalides is presented. This review includes 114 references to recent advanced studies. These data facilitated the confirmation of previously suggested trends in the molecular properties of the title compounds and a reliable estimation of the data of less-studied  $LnX_3$  molecules.